. Amendments to the Claims:

This listing of claims will replace all prior versions and listings of claims in the application:

Listing of Claims:

1. (Currently amended) A compounds compound of the formula:

or the pharmaceutically acceptable acid salts thereof wherein: $R_1 \ \text{is halogen or} \ C_1\text{-}C_4 \ \text{alkyl};$

- R_2 and R_3 are the same or different and represent hydrogen, halogen, $C_1\text{-}C_4$ alkyl, $C_1\text{-}C_4$ alkoxy, alkylthio, hydroxy, amino, monoalkylamino or dialkylamino,
 - with the proviso that R_2 and R_3 may not be 2-isopropoxyl and hydrogen respectively when R_1 is bromo;
- wherein in an assay for D2 receptor binding the compound exhibits a Ki value of greater than 300 nM.
- 2. (Original) A compound according to Claim 1, wherein $\ensuremath{R_1}$ is methyl.
 - 3. (Currently amended) A compound of the formula:

or the pharmaceutically acceptable salts thereof wherein $R_{\rm x}$ is fluorine, chlorine, bromine, or iodine; and

 R_2 and R_3 are the same or different and represent hydrogen, halogen, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, alkylthio, hydroxy, amino, monoalkylamino or dialkylamino;

wherein in an assay for D2 receptor binding the compound exhibits a Ki value of greater than 300 nM.

- 4. (Original) A compound according to Claim 3, wherein R_2 and R_3 may not be 2-isopropoxyl and hydrogen, respectively, when R_1 is bromo.
- 5. (Original) A compound according to claim 3, wherein $R_{\rm x}$ is chloride; R_2 and R_3 may not be 2-isopropoxyl and hydrogen, respectively, when R_1 is bromo; R_2 is chloride, methyl or methoxy; and R_3 is hydrogen or methyl.
- 6. (Original) A compound according to claim 5, wherein the phenyl group substituted with R_2 and R_3 is selected from the group consisting of:

7. (Currently amended) A compound of the formula:

$$R_{a} \xrightarrow{N} N \xrightarrow{\frac{11}{11}} R_{3}$$

or the pharmaceutically acceptable salts thereof wherein $R_a \mbox{ is } C_1\hbox{-} C_4 \mbox{ alkyl}; \mbox{ and}$

 R_2 and R_3 are the same or different and represent hydrogen, halogen, $C_1\text{-}C_4$ alkyl, $C_1\text{-}C_4$ alkoxy, alkylthio, hydroxy, amino, monoalkylamino or dialkylamino;

wherein in an assay for D2 receptor binding the compound exhibits a Ki value of greater than 300 nM.

- 8. (Original) A compound according to Claim 7, wherein $R_{\rm 1}$ is methyl.
- 9. (Original) A compound of according to Claim 7, wherein R_2 is chloride, fluoride, methyl or methoxy; and R_3 is hydrogen or methyl.
- 10. (Original) A compound according to claim 8, wherein the phenyl group substituted with R_2 and R_3 is selected from the group consisting of:

11. (Currently amended) A compound of the formula:

or the pharmaceutically acceptable salts thereof wherein:

 R_1 is C_1 - C_4 alkyl or halogen; and

wherein in an assay for D2 receptor binding the compound exhibits a Ki value of greater than 300 nM.

12. (Original) A compound according to Claim 11, wherein R_1 is chloro.

13-35. (Canceled)